Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1611bxv

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * * Welcome to STN International * * * * * * * * * *

NEWS	1		Web Page URLs for STN Seminar Schedule - N. America
NEWS	2		"Ask CAS" for self-help around the clock
NEWS	3	JUL 20	Powerful new interactive analysis and visualization software, STN AnaVist, now available
NEWS	4	AUG 11	Derwent World Patents Index(R) web-based training during August
NEWS	5	AUG 11	STN AnaVist workshops to be held in North America
NEWS	6	AUG 30	CA/CAplus -Increased access to 19th century research documents
NEWS	7	AUG 30	CASREACT - Enhanced with displayable reaction conditions
NEWS	8	SEP 09	ACD predicted properties enhanced in REGISTRY/ZREGISTRY

NEWS EXPRESS	JUNE 13 CURRENT WINDOWS VERSION IS V8.0, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 13 JUNE 2005
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
NEWS WWW	CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

* * * * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 16:39:47 ON 18 SEP 2005

=>
=>
=> FILE REG

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 16:40:13 ON 18 SEP 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 SEP 2005 HIGHEST RN 863378-74-9
DICTIONARY FILE UPDATES: 16 SEP 2005 HIGHEST RN 863378-74-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

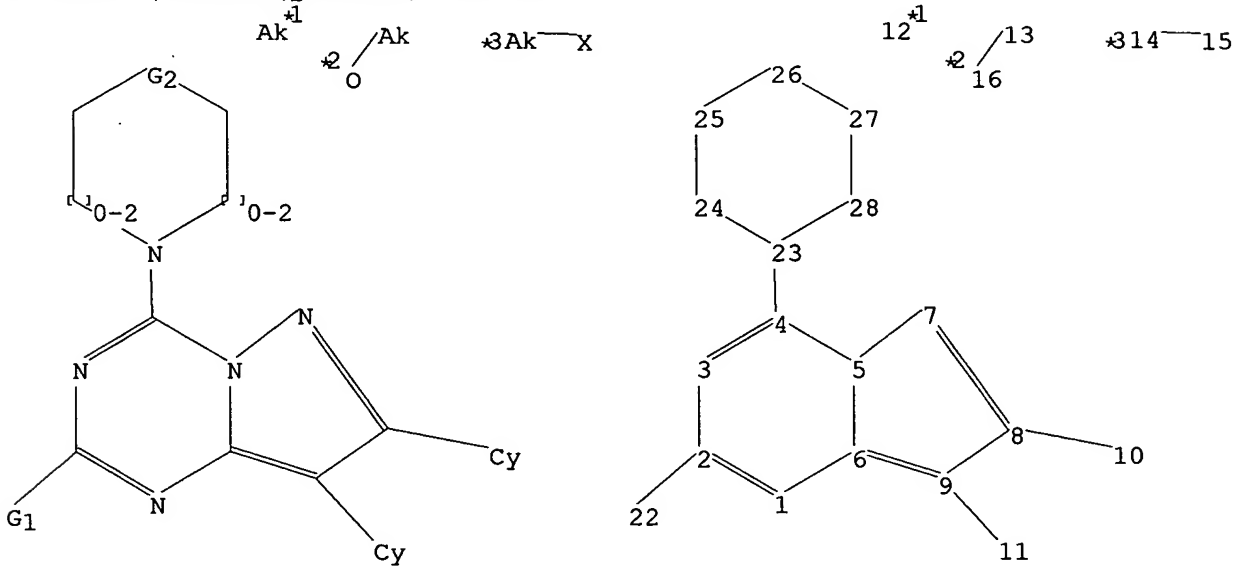
Please note that search-term pricing does apply when conducting SmartSELECT searches.

****
* *
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
* *
****

Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=>
Uploading C:\Documents and Settings\VBalasubramania\My Documents\STNEXP4\QUERIES\10763105.str



chain nodes :
10 11 12 13 14 15 16 22
ring nodes :
1 2 3 4 5 6 7 8 9 23 24 25 26 27 28
chain bonds :
2-22 4-23 8-10 9-11 13-16 14-15

ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 23-24 23-28 24-25 25-26 26-27
27-28
exact/norm bonds :
1-2 1-6 2-3 2-22 3-4 4-5 4-23 5-6 5-7 6-9 7-8 8-9 8-10 9-11 13-16
14-15 23-24 23-28 24-25 25-26 26-27 27-28

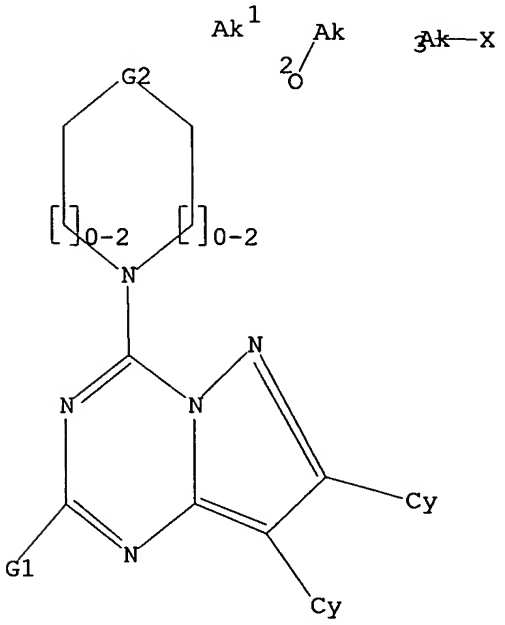
G1:H,[*1],[*2],[*3]

G2:C,O,S

Connectivity :
12:1 E exact RC ring/chain 13:1 E exact RC ring/chain 14:2 E exact RC ring/chain
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 22:CLASS 23:Atom
24:Atom 25:Atom 26:Atom 27:Atom 28:Atom
Generic attributes :
10:
Saturation : Unsaturated
11:
Saturation : Unsaturated

L1 STRUCTURE UPLOADED

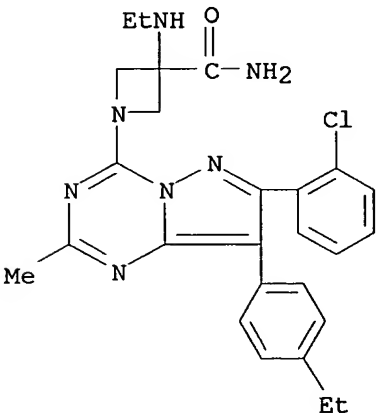
=> D L1
L1 HAS NO ANSWERS
L1 STR



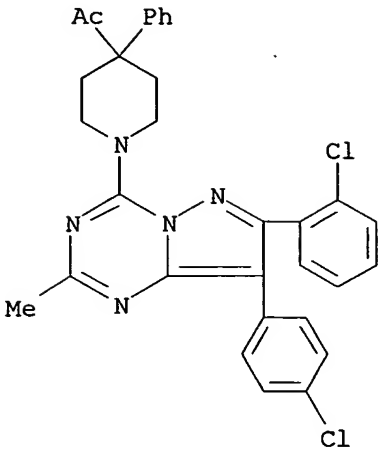
G1 H,[@1],[@2],[@3]
G2 C,O,S

Structure attributes must be viewed using STN Express query preparation.

```
=> S L1 SSS SAM
SAMPLE SEARCH INITIATED 16:40:39 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -       54 TO ITERATE

100.0% PROCESSED       54 ITERATIONS                                 2 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:          640 TO     1520
PROJECTED ANSWERS:               2 TO      124

L2            2 SEA SSS SAM L1

=> D SCAN

L2   2 ANSWERS   REGISTRY  COPYRIGHT 2005 ACS on STN
IN   3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-ethylphenyl)-2-
     methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI)
MF   C26 H28 Cl N7 O
```




```
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**


HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2   2 ANSWERS   REGISTRY  COPYRIGHT 2005 ACS on STN
IN   Ethanone, 1-[1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-
     a]-1,3,5-triazin-4-yl]-4-phenyl-4-piperidinyl]- (9CI)
MF   C31 H27 Cl2 N5 O
```



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> S L1 SSS FUL
FULL SEARCH INITIATED 16:40:54 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1082 TO ITERATE

100.0% PROCESSED 1082 ITERATIONS 49 ANSWERS
SEARCH TIME: 00.00.01

L3 49 SEA SSS FUL L1

=> FILE CAPLUS

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	161.33	161.54

FILE 'CAPLUS' ENTERED AT 16:41:02 ON 18 SEP 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1907 - 18 Sep 2005 VOL 143 ISS 13
FILE LAST UPDATED: 16 Sep 2005 (20050916/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> S L3
L4 2 L3

=> D L4 1-2 BIB HITSTR

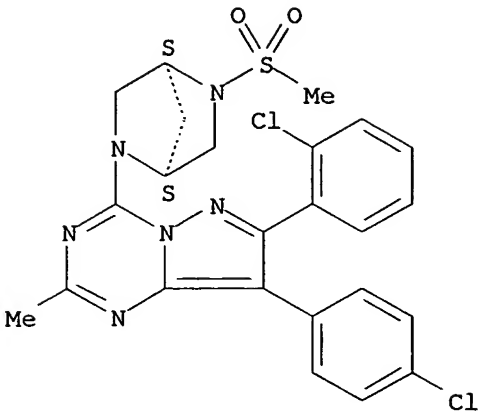
L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2005 ACS on STN
AN 2005:160837 CAPLUS
DN 142:233372
TI Pharmaceutical composition using a combination of an opioid receptor antagonist and a CB-1 receptor antagonist for the prevention and treatment of addiction in a mammal
IN Coe, Jotham Wadsworth; Iredale, Philip A.; McHardy, Stanton Furst; McLean, Stafford
PA Pfizer Inc, USA
SO U.S. Pat. Appl. Publ., 25 pp.
CODEN: USXXCO
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI US 2005043327	A1	20050224	US 2004-870209	20040617
WO 2005018645	A1	20050303	WO 2004-IB2596	20040809

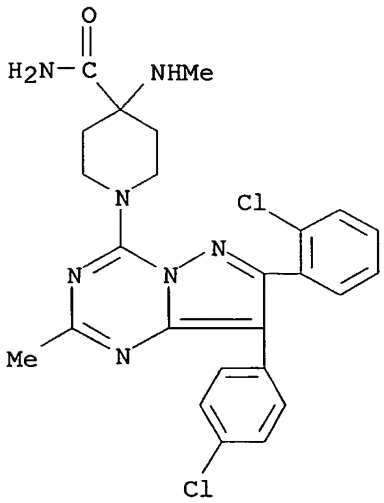
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

PRAI US 2003-496803P P 20030821
IT **736993-30-9 736993-50-3 736993-51-4**
736993-52-5 736993-53-6 736993-54-7
736993-55-8 736993-56-9 736993-57-0
736993-58-1 736993-63-8 736993-64-9
736993-65-0 736993-66-1 736993-67-2
736993-68-3 736993-69-4 736993-70-7
736994-31-3 845670-45-3
RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(opioid receptor antagonist-CB-1 receptor antagonist combination for prevention and treatment of addiction)
RN 736993-30-9 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-(methylsulfonyl)-, (1S,4S)-(9CI) (CA INDEX NAME)

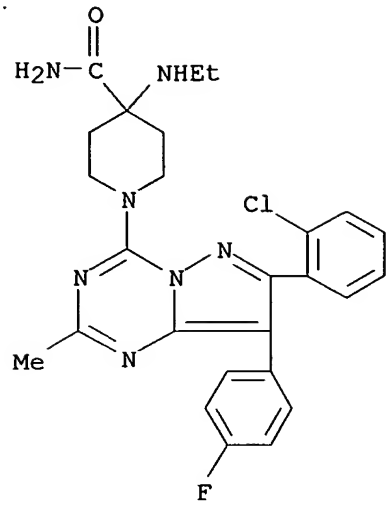
Absolute stereochemistry.



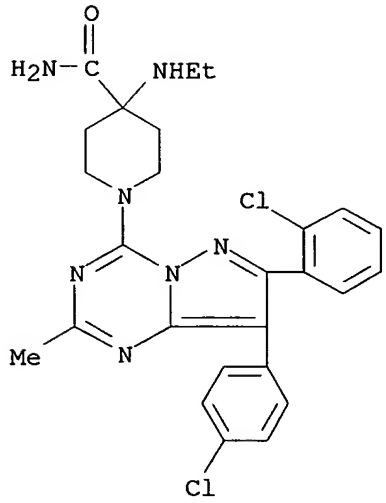
RN 736993-50-3 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(methylamino)- (9CI) (CA INDEX NAME)



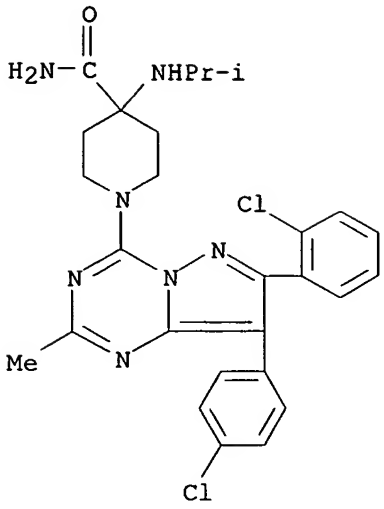
RN 736993-51-4 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-fluorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)- (9CI) (CA INDEX NAME)



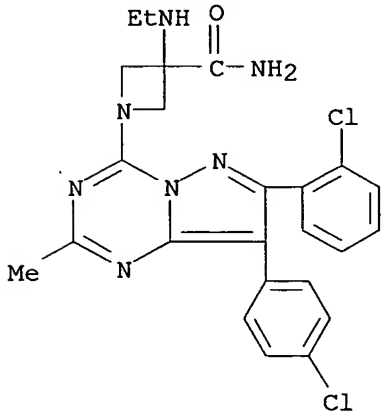
RN 736993-52-5 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)- (9CI) (CA INDEX NAME)



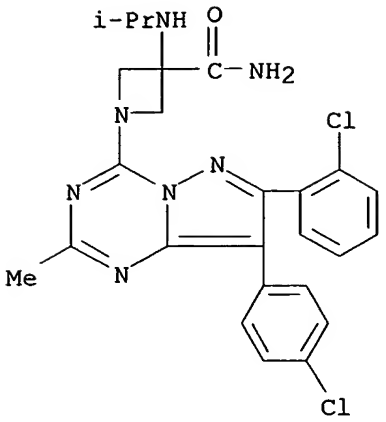
RN 736993-53-6 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-[(1-methylethyl)amino]- (9CI) (CA INDEX NAME)



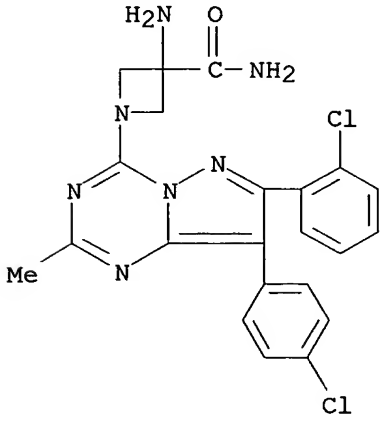
RN 736993-54-7 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



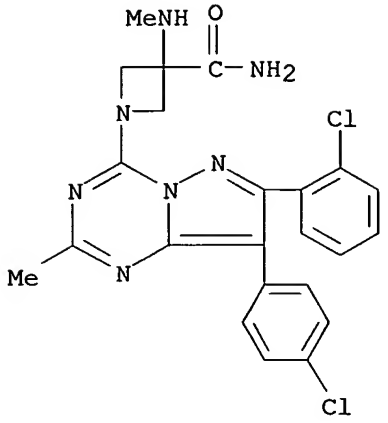
RN 736993-55-8 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-[(1-methylethyl)amino]- (9CI) (CA INDEX NAME)



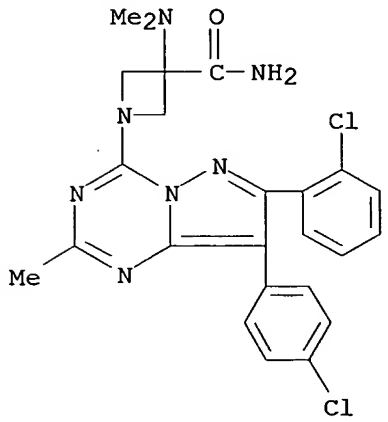
RN 736993-56-9 CAPLUS
CN 3-Azetidinecarboxamide, 3-amino-1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]- (9CI) (CA INDEX NAME)



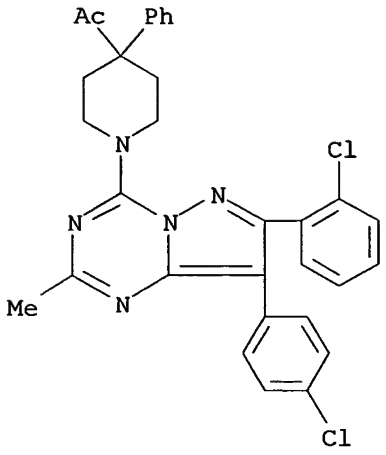
RN 736993-57-0 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(methylamino)- (9CI) (CA INDEX NAME)



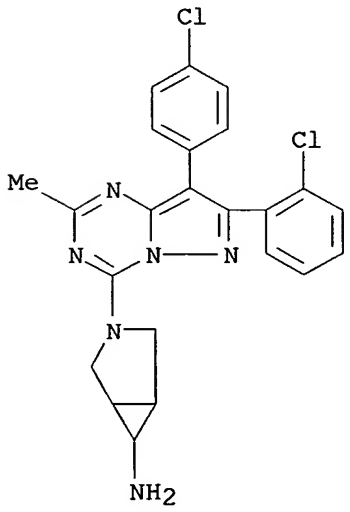
RN 736993-58-1 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(dimethylamino)- (9CI) (CA INDEX NAME)



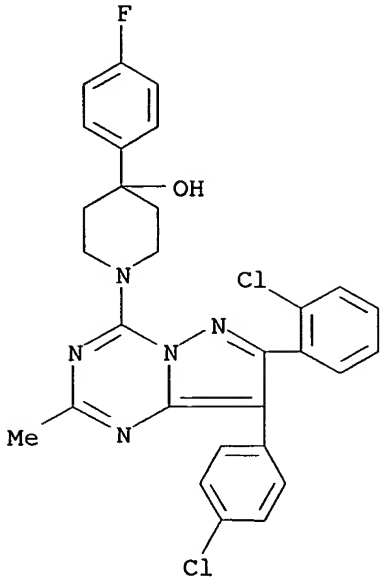
RN 736993-63-8 CAPLUS
CN Ethanone, 1-[1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-phenyl-4-piperidinyl]- (9CI) (CA INDEX NAME)



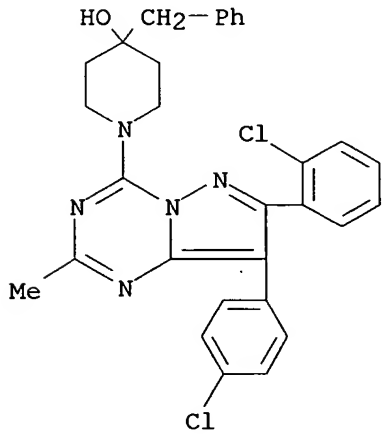
RN 736993-64-9 CAPLUS
CN 3-Azabicyclo[3.1.0]hexan-6-amine, 3-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]- (9CI) (CA INDEX NAME)



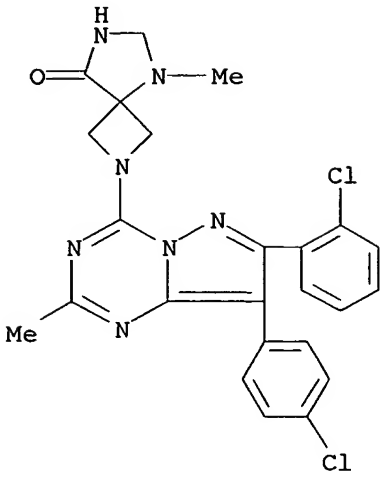
RN 736993-65-0 CAPLUS
CN 4-Piperidinol, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(4-fluorophenyl)- (9CI) (CA INDEX NAME)



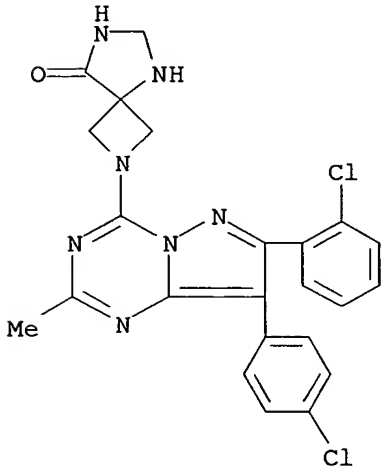
RN 736993-66-1 CAPLUS
CN 4-Piperidinol, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(phenylmethyl)- (9CI) (CA INDEX NAME)



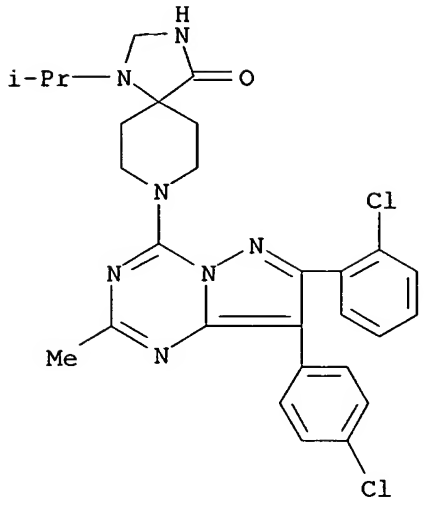
RN 736993-67-2 CAPLUS
CN 2,5,7-Triazaspiro[3.4]octan-8-one, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-methyl- (9CI) (CA INDEX NAME)



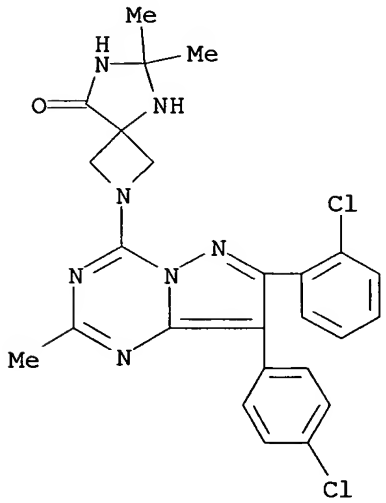
RN 736993-68-3 CAPLUS
CN 2,5,7-Triazaspiro[3.4]octan-8-one, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]- (9CI) (CA INDEX NAME)



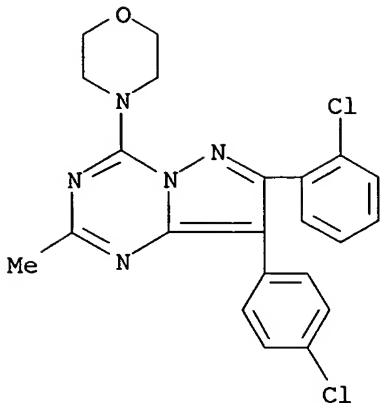
RN 736993-69-4 CAPLUS
CN 1,3,8-Triazaspiro[4.5]decan-4-one, 8-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



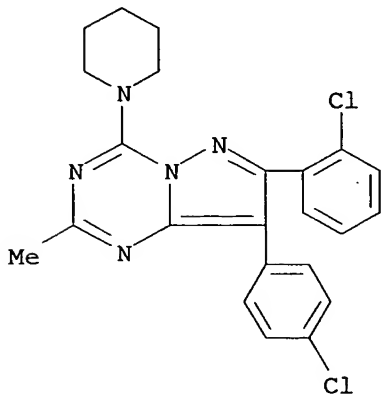
RN 736993-70-7 CAPLUS
CN 2,5,7-Triazaspiro[3.4]octan-8-one, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-6,6-dimethyl- (9CI) (CA INDEX NAME)



RN 736994-31-3 CAPLUS
CN Pyrazolo[1,5-a]-1,3,5-triazine, 7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methyl-4-(4-morpholinyl)- (9CI) (CA INDEX NAME)



RN 845670-45-3 CAPLUS
CN Pyrazolo[1,5-a]-1,3,5-triazine, 7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methyl-4-(1-piperidinyl)- (9CI) (CA INDEX NAME)



L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2005 ACS on STN
AN 2004:654773 CAPLUS
DN 141:190806
TI Preparation of pyrazolotriazines as cannabinoid CB1 receptor antagonists
IN Griffith, David A.
PA Pfizer Inc, USA
SO U.S. Pat. Appl. Publ., 72 pp.
CODEN: USXXCO
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2004157839	A1	20040812	US 2004-763105	20040121
	WO 2004069837	A1	20040819	WO 2004-IB269	20040126

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NL 1025404 A1 20040809 NL 2004-1025404 20040204
NL 1025404 C2 20050314
PRAI US 2003-445728P P 20030206
OS MARPAT 141:190806
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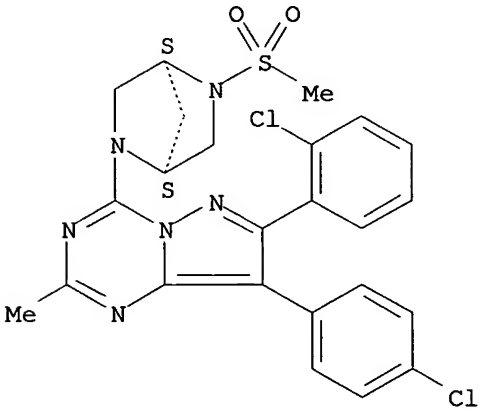
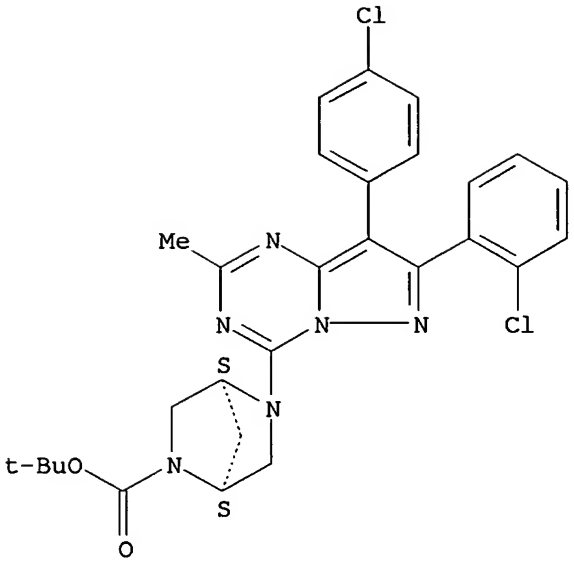
736993-62-7P 736993-63-8P 736993-64-9P
736993-65-0P 736993-66-1P 736993-67-2P
736993-68-3P 736993-69-4P 736993-70-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(claimed compound; preparation of pyrazolotriazines as cannabinoid CB1 receptor antagonists)
RN 736993-30-9 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-(methylsulfonyl)-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



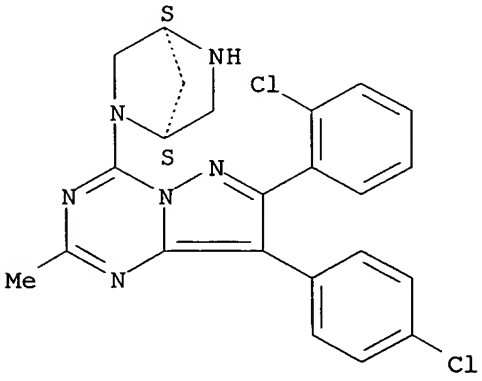
RN 736993-35-4 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 5-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-, 1,1-dimethylethyl ester, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



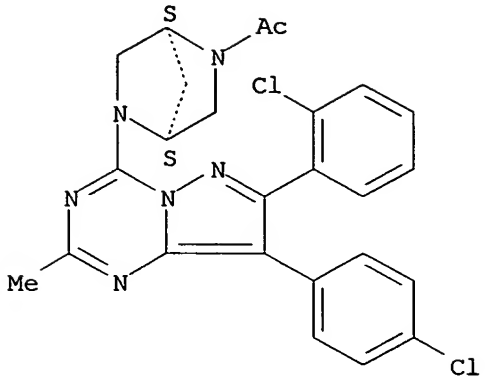
RN 736993-36-5 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



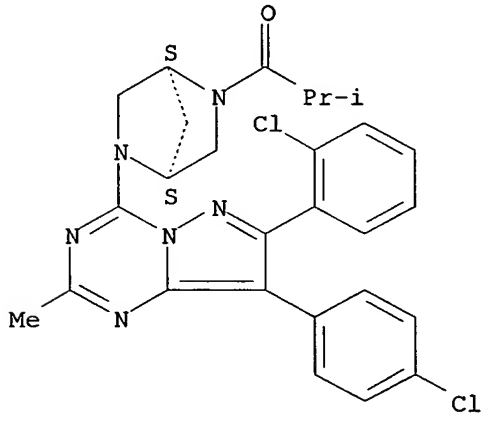
RN 736993-37-6 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-acetyl-5-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



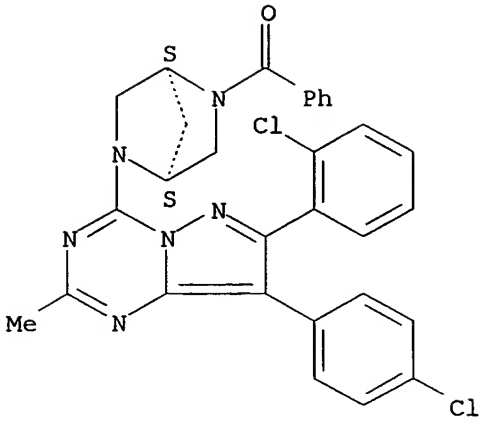
RN 736993-38-7 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-(2-methyl-1-oxopropyl)-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



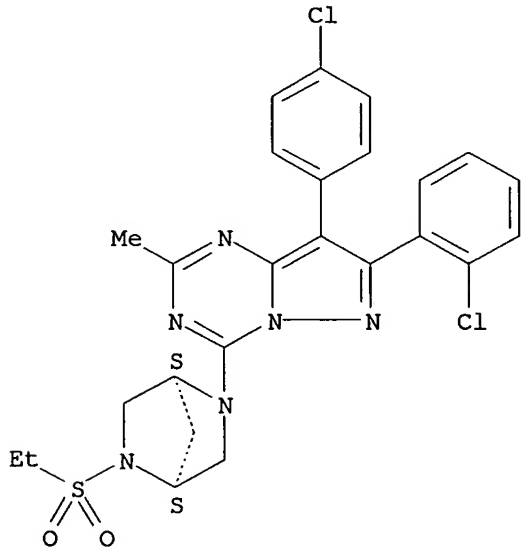
RN 736993-40-1 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-benzoyl-5-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



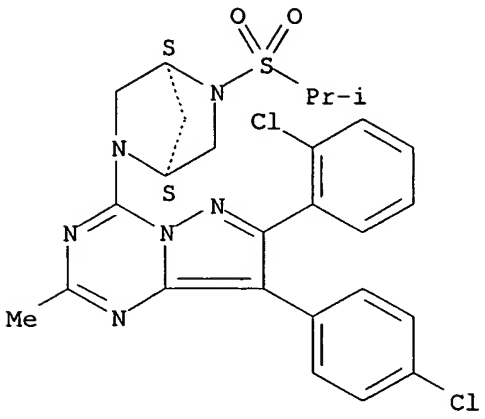
RN 736993-41-2 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-(ethylsulfonyl)-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



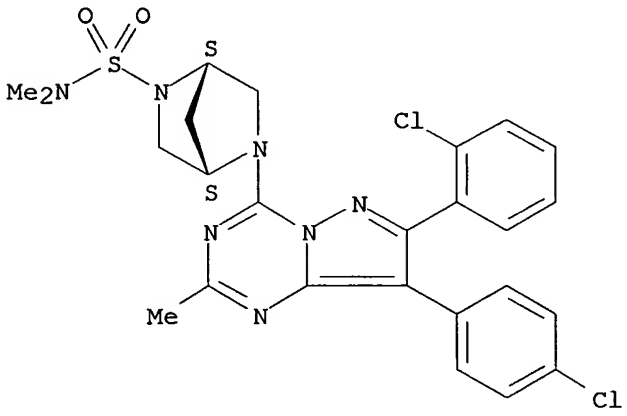
RN 736993-42-3 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-[(1-methylethyl)sulfonyl]-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

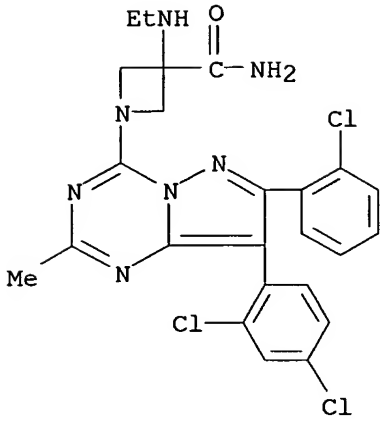


RN 736993-43-4 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane-2-sulfonamide, 5-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-N,N-dimethyl-, (1S,4S)- (9CI) (CA INDEX NAME)

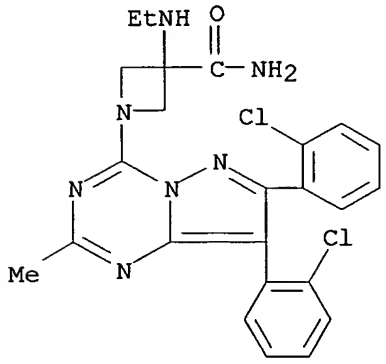
Absolute stereochemistry.



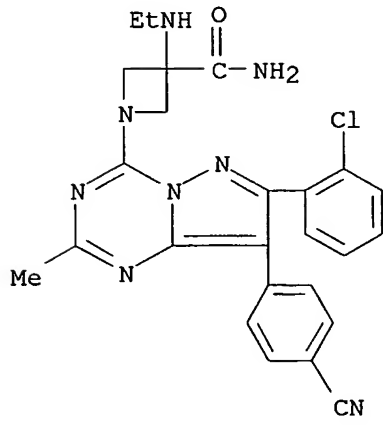
RN 736993-44-5 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(2,4-dichlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



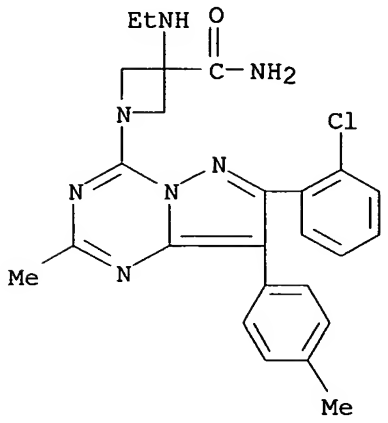
RN 736993-45-6 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7,8-bis(2-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



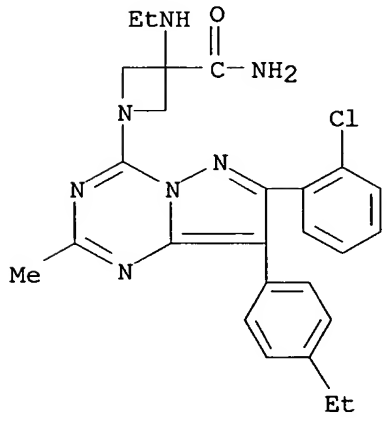
RN 736993-46-7 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-cyanophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



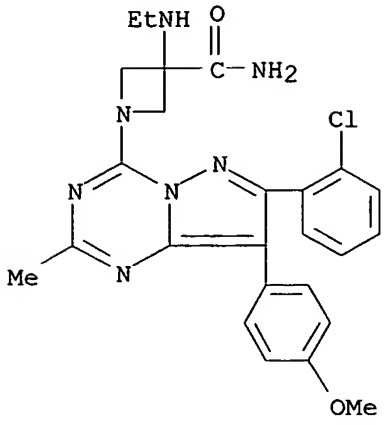
RN 736993-47-8 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-2-methyl-8-(4-methylphenyl)pyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



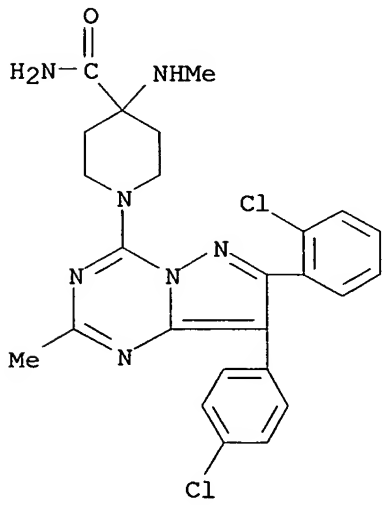
RN 736993-48-9 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-ethylphenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



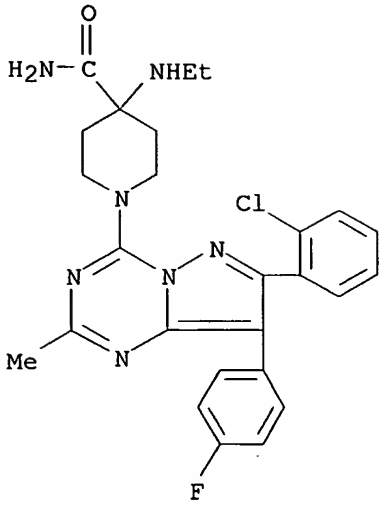
RN 736993-49-0 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-methoxyphenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



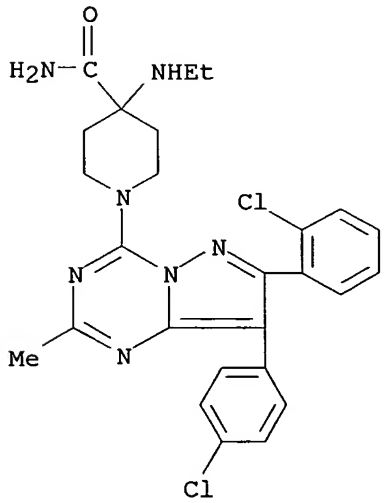
RN 736993-50-3 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(methylamino)- (9CI) (CA INDEX NAME)



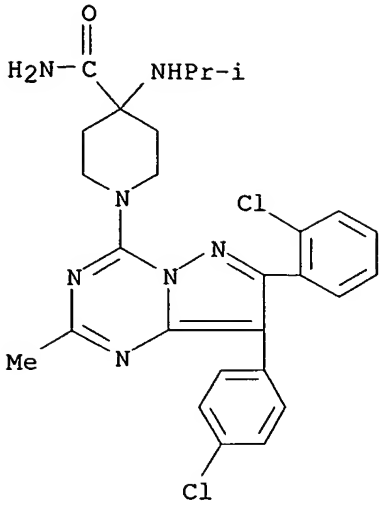
RN 736993-51-4 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-fluorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)- (9CI) (CA INDEX NAME)



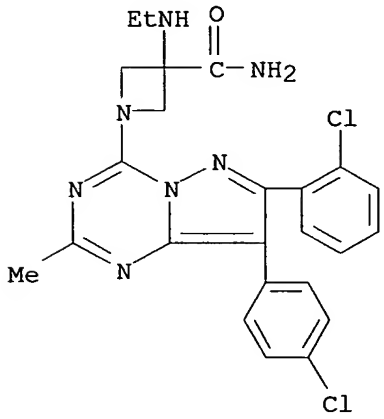
RN 736993-52-5 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)- (9CI) (CA INDEX NAME)



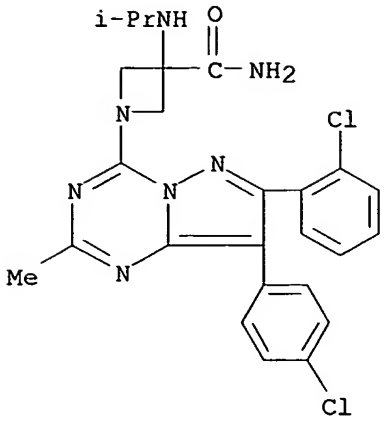
RN 736993-53-6 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-[(1-methylethyl)amino]- (9CI) (CA INDEX NAME)



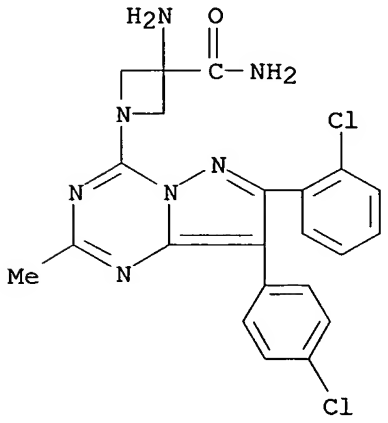
RN 736993-54-7 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



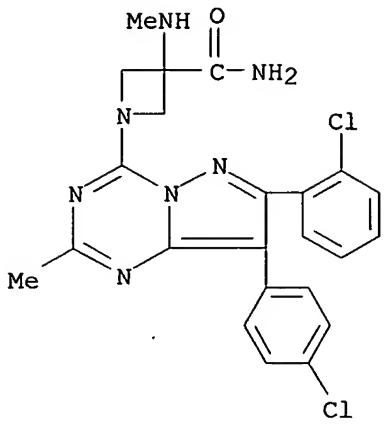
RN 736993-55-8 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-[(1-methylethyl)amino]- (9CI) (CA INDEX NAME)



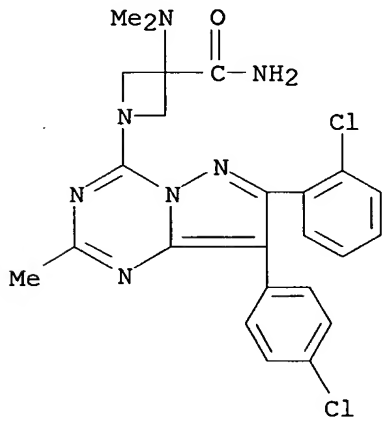
RN 736993-56-9 CAPLUS
CN 3-Azetidinecarboxamide, 3-amino-1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]- (9CI) (CA INDEX NAME)



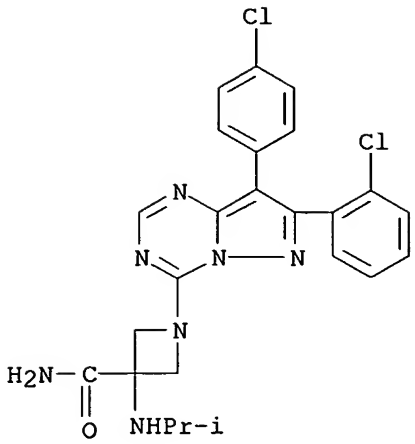
RN 736993-57-0 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(methylamino)- (9CI) (CA INDEX NAME)



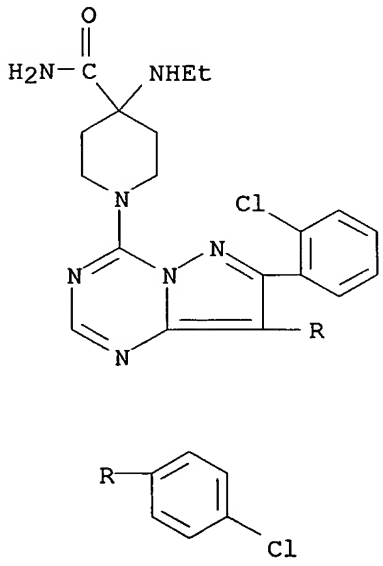
RN 736993-58-1 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(dimethylamino)- (9CI) (CA INDEX NAME)



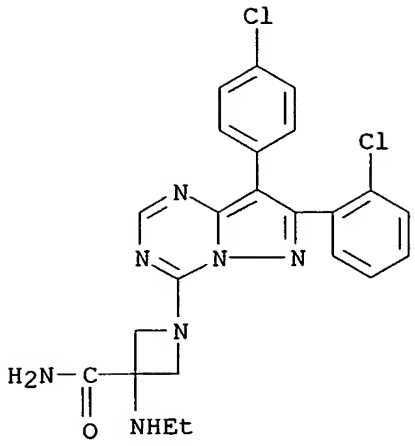
RN 736993-59-2 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)pyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-[(1-methylethyl)amino]- (9CI) (CA INDEX NAME)



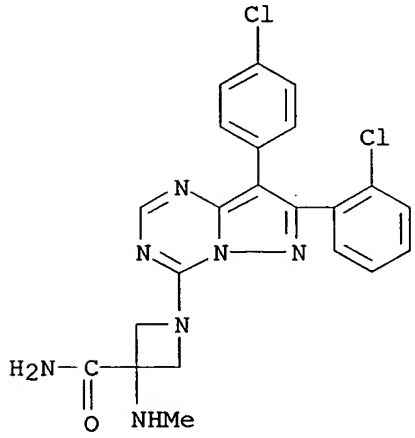
RN 736993-60-5 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)pyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)- (9CI) (CA INDEX NAME)



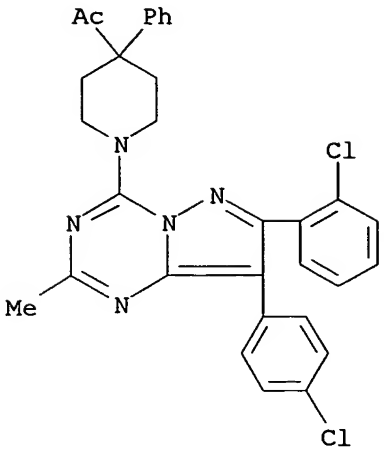
RN 736993-61-6 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)pyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)



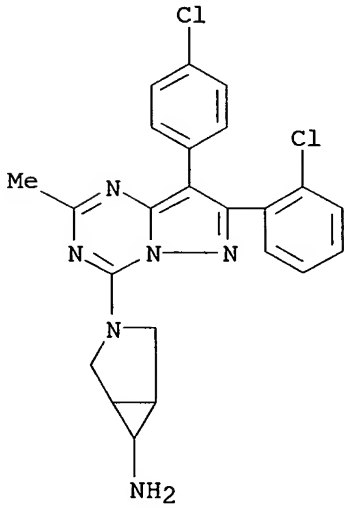
RN 736993-62-7 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)pyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(methylamino)- (9CI) (CA INDEX NAME)



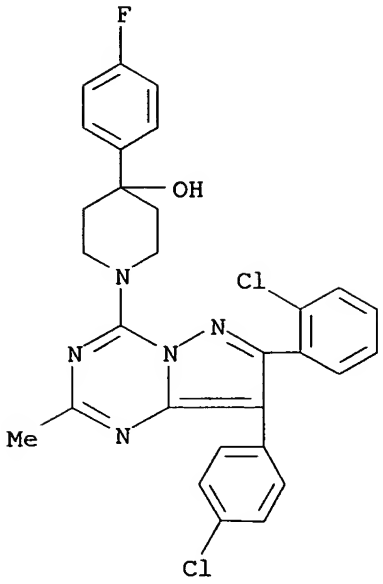
RN 736993-63-8 CAPLUS
CN Ethanone, 1-[1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-phenyl-4-piperidinyl]- (9CI) (CA INDEX NAME)



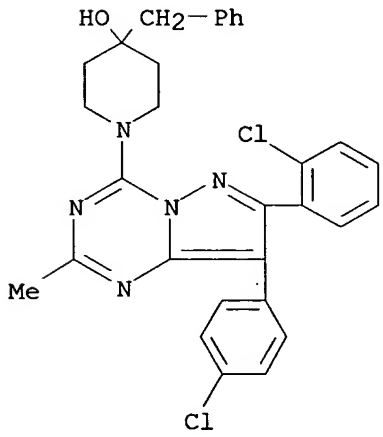
RN 736993-64-9 CAPLUS
CN 3-Azabicyclo[3.1.0]hexan-6-amine, 3-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]- (9CI) (CA INDEX NAME)



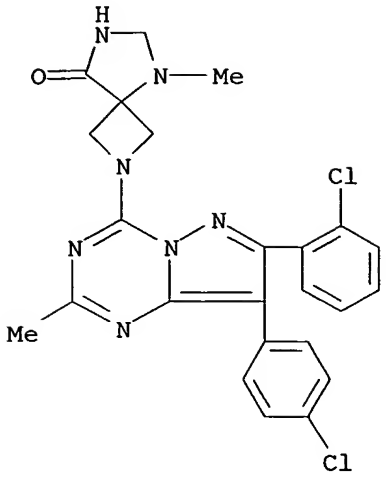
RN 736993-65-0 CAPLUS
CN 4-Piperidinol, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(4-fluorophenyl)- (9CI) (CA INDEX NAME)



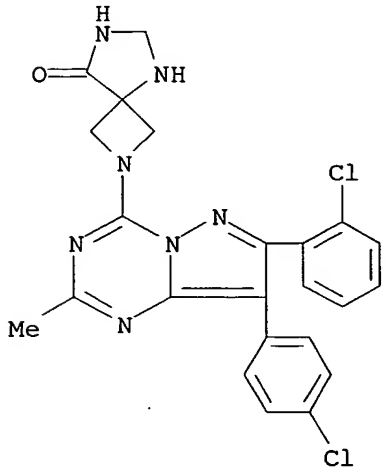
RN 736993-66-1 CAPLUS
CN 4-Piperidinol, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(phenylmethyl)- (9CI) (CA INDEX NAME)



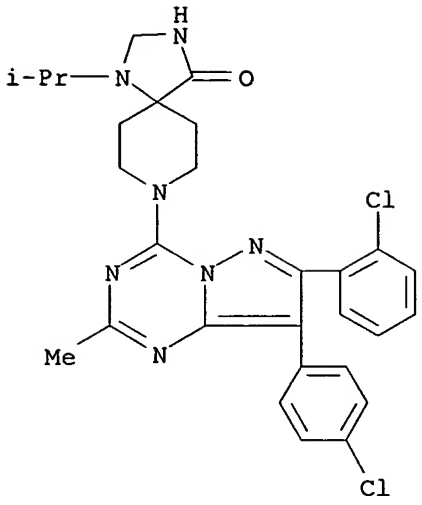
RN 736993-67-2 CAPLUS
CN 2,5,7-Triazaspiro[3.4]octan-8-one, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-methyl- (9CI) (CA INDEX NAME)



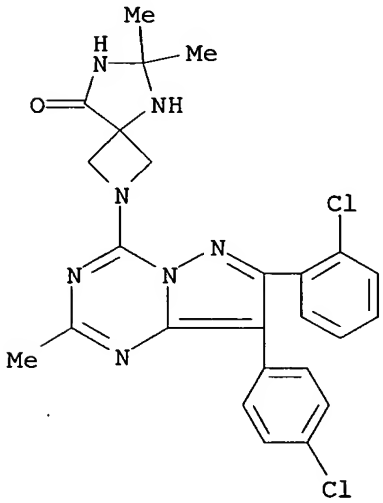
RN 736993-68-3 CAPLUS
CN 2,5,7-Triazaspiro[3.4]octan-8-one, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]- (9CI) (CA INDEX NAME)



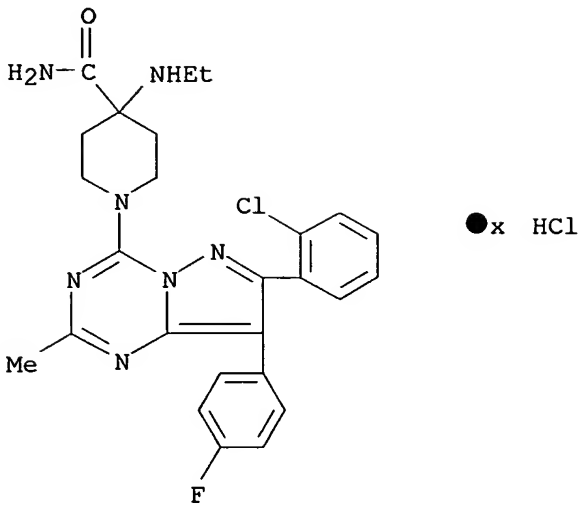
RN 736993-69-4 CAPLUS
CN 1,3,8-Triazaspiro[4.5]decan-4-one, 8-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



RN 736993-70-7 CAPLUS
CN 2,5,7-Triazaspiro[3.4]octan-8-one, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-6,6-dimethyl- (9CI) (CA INDEX NAME)

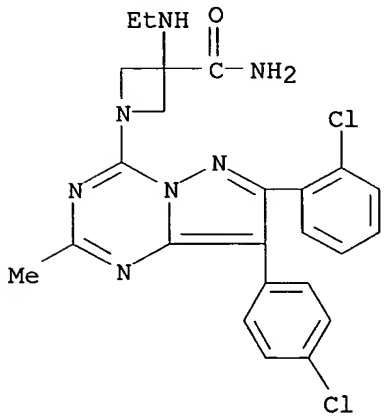


IT **736994-23-3P 736994-24-4P 736994-25-5P**
736994-26-6P 736994-27-7P 736994-29-9P
736994-30-2P 736994-31-3P 736994-32-4P
736994-33-5P 736994-34-6P 736994-35-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of pyrazolotriazines as cannabinoid CB1 receptor antagonists)
RN 736994-23-3 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-fluorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)-, hydrochloride (9CI) (CA INDEX NAME)



● x HCl

RN 736994-24-4 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)-, monohydrochloride (9CI) (CA INDEX NAME)

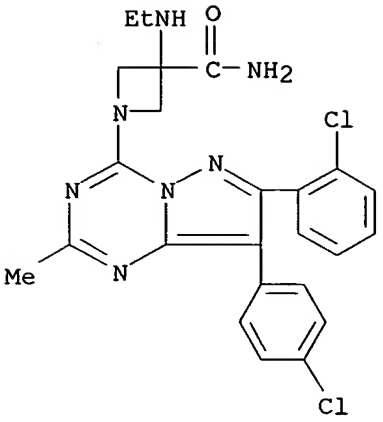


● HCl

RN 736994-25-5 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)-, benzenesulfonate (9CI) (CA INDEX NAME)

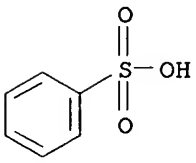
CM 1

CRN 736993-54-7
CMF C24 H23 Cl2 N7 O



CM 2

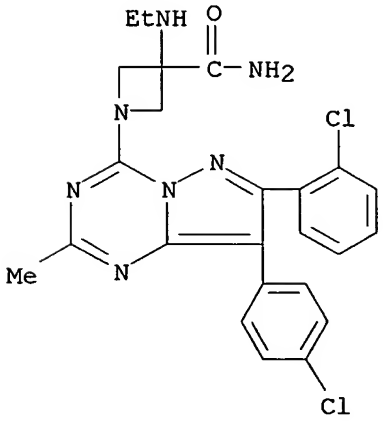
CRN 98-11-3
CMF C6 H6 O3 S



RN 736994-26-6 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)-, monomethanesulfonate (9CI) (CA INDEX NAME)

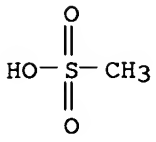
CM 1

CRN 736993-54-7
CMF C24 H23 Cl2 N7 O

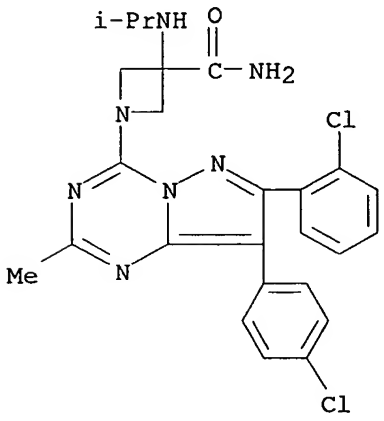


CM 2

CRN 75-75-2
CMF C H4 O3 S

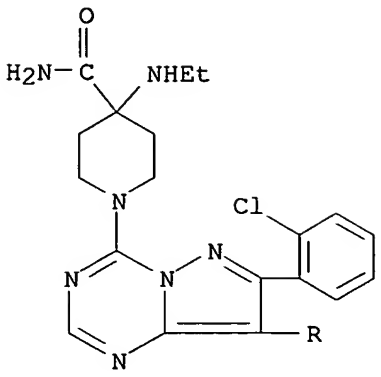


RN 736994-27-7 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-[(1-methylethyl)amino]-, monohydrochloride (9CI) (CA INDEX NAME)

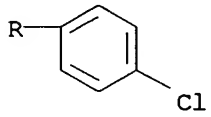


● HCl

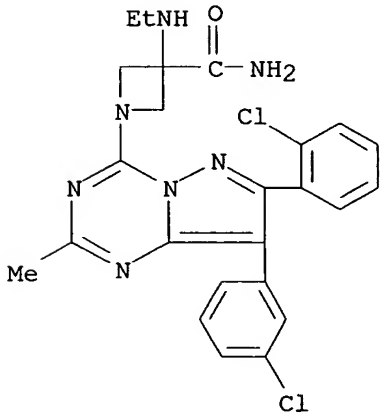
RN 736994-29-9 CAPLUS
CN 4-Piperidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(4-chlorophenyl)pyrazolo[1,5-a]-1,3,5-triazin-4-yl]-4-(ethylamino)-, hydrochloride (9CI) (CA INDEX NAME)



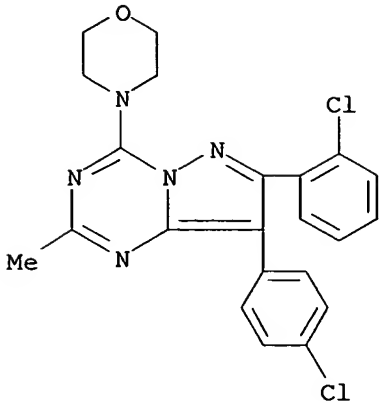
● x HCl



RN 736994-30-2 CAPLUS
CN 3-Azetidinecarboxamide, 1-[7-(2-chlorophenyl)-8-(3-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-3-(ethylamino)- (9CI) (CA INDEX NAME)

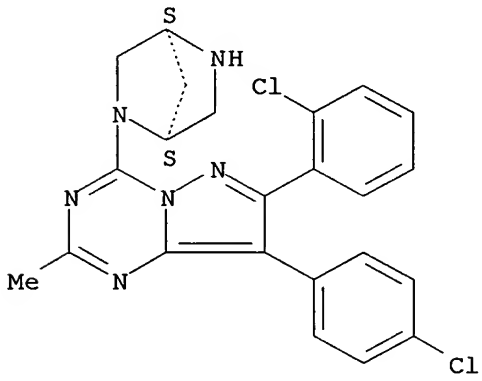


RN 736994-31-3 CAPLUS
CN Pyrazolo[1,5-a]-1,3,5-triazine, 7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methyl-4-(4-morpholinyl)- (9CI) (CA INDEX NAME)



RN 736994-32-4 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-, dihydrochloride, (1S,4S)- (9CI) (CA INDEX NAME)

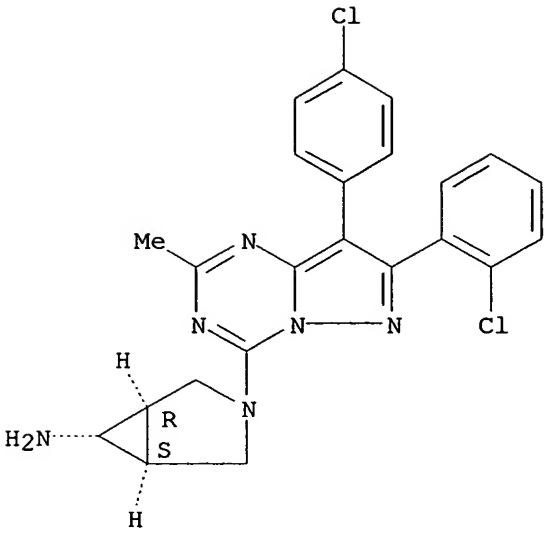
Absolute stereochemistry.



●2 HCl

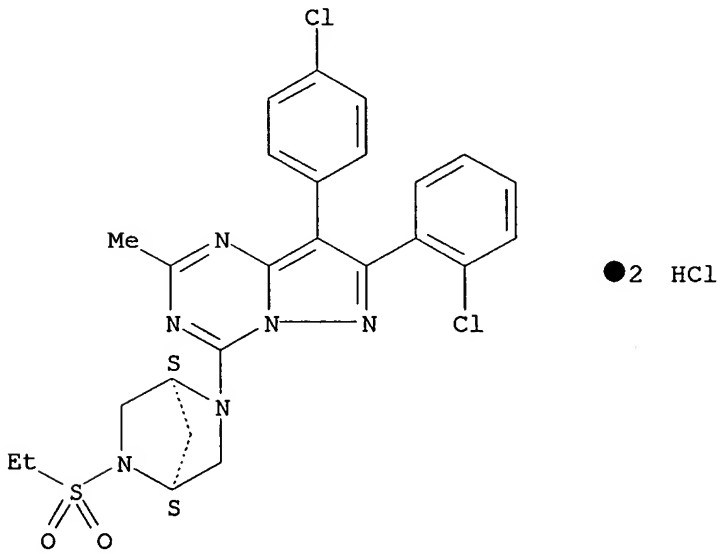
RN 736994-33-5 CAPLUS
CN 3-Azabicyclo[3.1.0]hexan-6-amine, 3-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-, (1α,5α,6α)- (9CI) (CA INDEX NAME)

Relative stereochemistry.



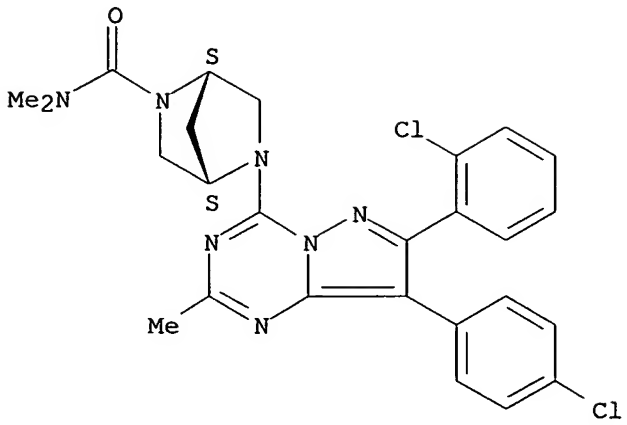
RN 736994-34-6 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane, 2-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-5-(ethylsulfonyl)-, dihydrochloride, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 736994-35-7 CAPLUS
CN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxamide, 5-[7-(2-chlorophenyl)-8-(4-chlorophenyl)-2-methylpyrazolo[1,5-a]-1,3,5-triazin-4-yl]-N,N-dimethyl-, (1S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



=> LOG Y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	7.23	168.77

STN INTERNATIONAL LOGOFF AT 16:41:21 ON 18 SEP 2005

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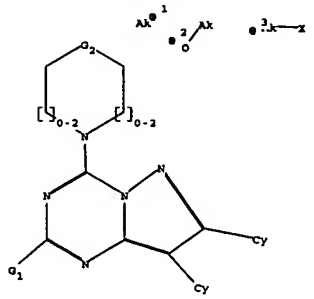

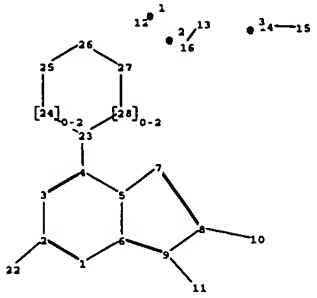

chain nodes :
10 11 12 13 14 15 16 22
ring nodes :
1 2 3 4 5 6 7 8 9 23 24 25 26 27 28
chain bonds :
2-22 4-23 8-10 9-11 13-16 14-15
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 23-24 23-28 24-25 25-26 26-27 27-28
exact/norm bonds :
1-2 1-6 2-3 2-22 3-4 4-5 4-23 5-6 5-7 6-9 7-8 8-9 8-10 9-11 13-16 14-15
23-24 23-28 24-25 25-26 26-27 27-28

G1:H,[*1],[*2],[*3]

G2:C,O,S

Connectivity :
12:1 E exact RC ring/chain 13:1 E exact RC ring/chain 14:2 E exact RC ring/chain
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 22:CLASS 23:Atom 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom
Generic attributes :
10:
Saturation : Unsaturated
11:
Saturation : Unsaturated

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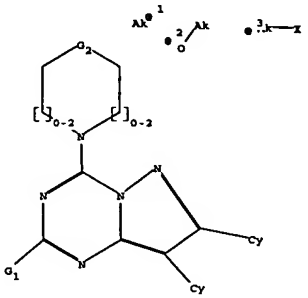

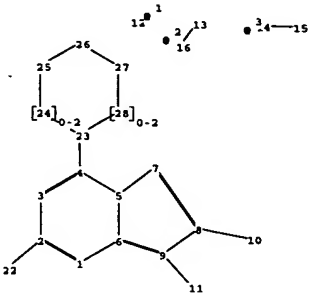

chain nodes :
10 11 12 13 14 15 16 22
ring nodes :
1 2 3 4 5 6 7 8 9 23 24 25 26 27 28
chain bonds :
2-22 4-23 8-10 9-11 13-16 14-15
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 23-24 23-28 24-25 25-26 26-27 27-28
exact/norm bonds :
1-2 1-6 2-3 2-22 3-4 4-5 4-23 5-6 5-7 6-9 7-8 8-9 8-10 9-11 13-16 14-15
23-24 23-28 24-25 25-26 26-27 27-28

G1:H,[*1],[*2],[*3]

G2:C,O,S

Connectivity :
12:1 E exact RC ring/chain 13:1 E exact RC ring/chain 14:2 E exact RC ring/chain
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 22:CLASS 23:Atom 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom
Generic attributes :
10:
Saturation : Unsaturated
11:
Saturation : Unsaturated